

Non-BCS Temperature Dependence of Energy Gap in Thin Film Electron-Doped Cuprates

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Abstract We investigate the dependence of the energy gap (G) on the temperature (T) for the electron-doped high-temperature superconductors. The following compounds, in the form of the thin films, have been taken into consideration: $\text{La}_{2-x}\text{Ce}_x\text{CuO}_4$ (LCCO), $\text{Pr}_{2-x}\text{Ce}_x\text{CuO}_4$ (PCCO), and $\text{Nd}_{2-x}\text{Ce}_x\text{CuO}_4$ (NCCO). It was found that $G(T)$ deviates from the BCS prediction more, if a concentration of cerium assumes the lower values. For the lowest concentration (in the case of LCCO and NCCO), the function $G(T)$ is not quite like the BCS curve, which is connected with the existence of the residual Nernst region. Next, it has been pointed out that the NCCO superconductor becomes structurally unstable for the maximum concentration of cerium, which is leading to the anomalous dependence of the energy gap on the temperature and the induction of the wide Nernst region.

Keywords Superconductivity · Electron-doped cuprates · Thin films · Thermodynamic properties

1 Introduction

The high-temperature superconductivity in cuprates has been discovered in 1986 [1, 2]. Cuprates can be divided into two groups: the hole-doped and the electron-doped superconductors [3–5]. Both families have distinctly different phase diagrams [6].

In the case of the hole-doped superconductors, the narrow antiferromagnetic region can be observed for the low values of the doping (δ). With the increase of δ , the distinct superconducting phase induces for the lower temperatures. The Cooper condensate borders from the high values of T with the Nernst region [7, 8]. Above, the pseudogap is observed [9–12]. Both the Nernst and the pseudogap region exists at least to the value of δ_{max} , which corresponding to the optimal doping.

On the other hand, the antiferromagnetic phase is very extended in the group of the electron-doped superconductors, and it is bordered with the relatively small region corresponding to the superconducting state [13]. The Nernst region is observed above the superconducting dome [14]. However, there are no clear results indicating the existence of the pseudogap [13].

It should be noted that the hole-doped superconductors characterize the much higher values of the critical temperature (T_C) than the electron-doped compounds. Currently, the highest T_C equal to ~ 160 K has been measured for $\text{HgBa}_2\text{Ca}_2\text{Cu}_3\text{O}_{8+y}$, located under the pressure of ~ 31 GPa [15].

In the family of the electron-doped superconductors, the highest critical temperature is equal to ~ 30 K, and it has been obtained for $\text{La}_{2-x}\text{Ce}_x\text{CuO}_4$ (LCCO) [16, 17]. However, $\text{Sr}_{0.9}\text{La}_{0.1}\text{CuO}_2$ possesses $T_C \sim 40$ K, but the synthesis of this compound is extremely difficult [18].

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The hole-doped superconductors have the d -wave symmetry of the order parameter [19], while the symmetry in the electron-doped compounds is still under debate. In particular, the results obtained with the help of the penetration depth [20–23], the tunneling spectroscopy [24, 25], and the Raman scattering [26] point at the s -wave symmetry. On the other hand, the other data (the penetration depth [27, 28], ARPES [29–31], the tricrystal experiment [32], and the Raman scattering [33]) emphasize the importance of the d -wave symmetry.

In the case of cuprates, it was very quickly realized that the thermodynamic properties of the superconducting phase cannot be accurately calculated in the framework of the BCS theory or the Eliashberg formalism [34–39]. The above fact results from the exotic pairing mechanism and may be explained in two ways. The first theory emphasizes the independent role of the strong electron correlations, which are usually modeled by the Hubbard Hamiltonian [6, 40, 41]. In particular, the performed *ab initio* calculations show that in the case of the single copper-oxygen plane the hopping integral between the nearest neighbours (t_0) is equal to ~ 400 meV, and the on-site Coulomb energy (U_0) has the value of ~ 5 eV [42]. Note that this reasoning naturally explains the existence of the antiferromagnetic phase in cuprates. In particular, for the half-filled electron band, the Hubbard Hamiltonian can be reduced to the Heisenberg operator with the antiferromagnetic hopping integral ($J_0 \equiv 2t_0^2/U_0$) [6, 43].

Unfortunately, under this scheme, it has been failed to describe the superconducting phase in the consistent way [44–46]. For example, there are great difficulties in explaining the energy gap dependence on the temperature for the

lower values of δ [39, 47–50]. The issue, which is not settled, is also the origin of the pseudogap.

The second path of research is based on the experimental results, which underline the importance of the electron-phonon interaction [51–57]. Note that in the present case, the Coulomb energy should be treated as the source of the depairing correlations.

In order to substantiate the discussion, let us consider the simplest version of the Hubbard model:

$$H_0 = \varepsilon_0 \sum_{i\sigma} n_{i\sigma} + t_0 \sum_{ij\sigma} c_{i\sigma}^\dagger c_{j\sigma} + U_0 \sum_i n_{i\uparrow} n_{i\downarrow}, \quad (1)$$

where $c_{i\sigma}$ and $c_{i\sigma}^\dagger$ denote the annihilation and creation operators for the electron state. The index i labels the spatial lattice site and σ is the spin. Additionally: $n_{i\sigma} \equiv c_{i\sigma}^\dagger c_{i\sigma}$. The symbol ε_0 represents the reference energy.

The parameters included in the Hamiltonian (1) should be calculated by using the formulas:

$$\varepsilon_0 \equiv \int d^3\mathbf{r} \Phi_i^*(\mathbf{r}) \left[-\frac{1}{2} \nabla_{\mathbf{r}}^2 + V(\mathbf{r}) \right] \Phi_i(\mathbf{r}), \quad (2)$$

$$t_0 \equiv \int d^3\mathbf{r} \Phi_i^*(\mathbf{r}) \left[-\frac{1}{2} \nabla_{\mathbf{r}}^2 + V(\mathbf{r}) \right] \Phi_j(\mathbf{r}), \quad (3)$$

$$U_0 \equiv \int \int d^3\mathbf{r}_1 d^3\mathbf{r}_2 |\Phi_i(\mathbf{r}_1)|^2 V(\mathbf{r}_1 - \mathbf{r}_2) |\Phi_i(\mathbf{r}_2)|^2, \quad (4)$$

where the standard notation has been used as well as the Born-Oppenheimer approximation [58].

The Wannier function $\Phi_i(\mathbf{r})$ depends on the instantaneous position of the atom. Therefore, having regard to

Fig. 1 The influence of the cerium concentration on the value of **a** the critical temperature and **b** the energy gap for the superconductor LCCO [17, 67]

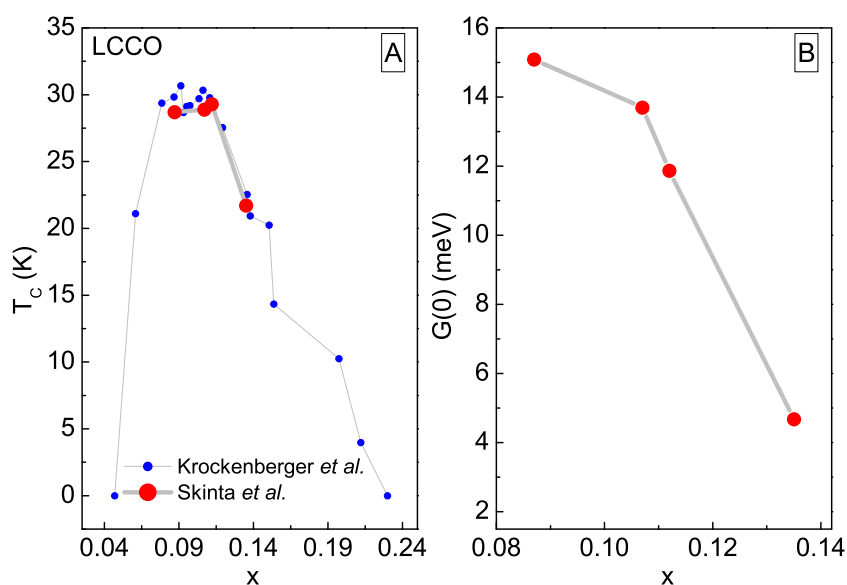


Table 1 The thermodynamic parameters of the LCCO superconductor. Additionally, it has been defined: $U^{(0)} \equiv \frac{U}{6} |\Delta(0)|^2$

x	T_C (K)	G(0) (meV)	V (meV)	$U^{(0)}$ (meV)	T^{**} (K)
0.087	28.7	15.1	436.6	123.9	28.9
0.107	28.9	13.7	438.0	95.6	—
0.112	29.3	11.9	440.7	56.7	—
0.135	21.7	4.7	387.8	−50.8	—

the lattice vibrations, the Hubbard Hamiltonian has to be generalized as follows: $H \rightarrow H_0 + \delta H$, where:

$$\delta H = g_{\varepsilon_0} \sum_{i\sigma} n_{i\sigma} \phi_i + g_{t_0} \sum_{ij\sigma} c_{i\sigma}^\dagger c_{j\sigma} \phi_i + g_{U_0} \sum_i n_{i\uparrow} n_{i\downarrow} \phi_i + \omega_0 \sum_i b_i^\dagger b_i. \quad (5)$$

The symbols g_{ε_0} , g_{t_0} , and g_{U_0} denote the electron-phonon coupling constants. Additionally: $\phi_i \equiv b_i^\dagger + b_i$, where b_i (b_i^\dagger) is the annihilation (creation) operator of the phonon state. The parameter ω_0 represents the value of the maximum phonon frequency.

Next, the Hamiltonian H can be rewritten in the momentum representation. The canonical transformation

eliminating the phonon degrees of freedom allows to obtain the fundamental thermodynamic equation [47]:

$$1 = \left(V + \frac{U}{6} |\Delta(T)|^2 \right) \int_{-\omega_0}^{\omega_0} d\varepsilon \rho(\varepsilon) \frac{\tanh\left(\frac{\beta}{2} \sqrt{\varepsilon^2 + E^2}\right)}{2\sqrt{\varepsilon^2 + E^2}}, \quad (6)$$

where $E \equiv \left(V + \frac{U}{6} |\Delta(T)|^2 \right) |\Delta(T)|$. The parameters V and U denote the electron-phonon potential (EPH) and the electron-electron-phonon potential (EPh), respectively: $V \equiv V_{\varepsilon_0} + V_{t_0}$ and $U \equiv V_{U_0} - U_0$.

The order parameter is defined by the formula: $\Delta \equiv \sum_{\mathbf{k}} \langle c_{-\mathbf{k}\downarrow} c_{\mathbf{k}\uparrow} \rangle$, while $c_{\mathbf{k}\sigma}$ denotes the annihilation operator for the electron state with the momentum \mathbf{k} . The symbol $\langle \dots \rangle$ represents the thermodynamic average [59, 60]. The summation over \mathbf{k} should be performed taking under consideration the condition: $\varepsilon_{\mathbf{k}} \leq |\omega_0|$, where $\varepsilon_{\mathbf{k}}$ is the electron band energy. The inverse temperature has been defined with

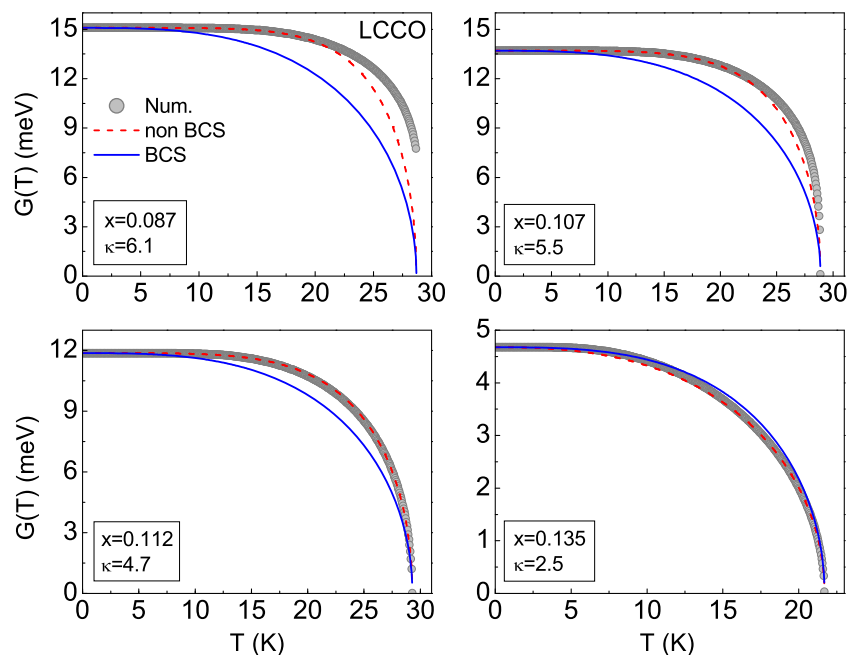
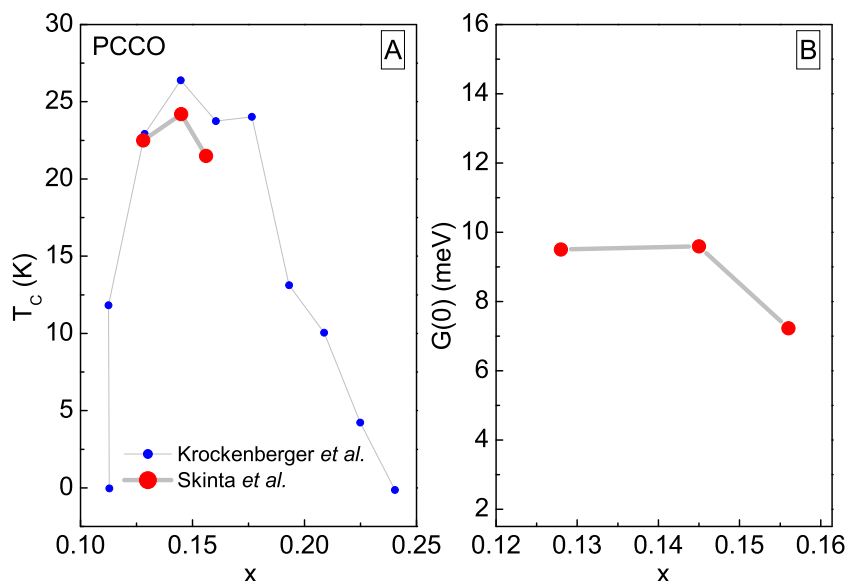
Fig. 2 The dependence of the LCCO energy gap on the temperature for the selected values of the Ce concentration

Fig. 3 The influence of the cerium concentration on the value of **a** the critical temperature and **b** the energy gap for the PCCO superconductor [17, 67]



the help of the expression: $\beta \equiv 1/(k_B T)$ where k_B is the Boltzmann constant.

Let us notice that the band energy for the two-dimensional square lattice can be written as $\varepsilon_{\mathbf{k}} = -2t_0 [\cos(k_x) + \cos(k_y)]$. Thus, the electron density of states has the form: $\rho(\varepsilon) = b_1 \ln|\varepsilon/b_2|$, where $b_1 \equiv -0.04687t_0^{-1}$ and $b_2 \equiv 21.17796t_0$ [61–65].

The presented model can explain (at the qualitative level) the essential properties of the high-temperature superconducting state [47]. Namely, the high critical temperature is related to the existence of the van Hove singularity in the electron density of states. The van Hove singularity is also responsible for the relatively low isotope coefficient. The anomalously high value of the energy gap ($G(0)$), where $G(0) \equiv 2(V + \frac{U}{6}|\Delta(0)|^2)|\Delta(0)|$ in respect to the critical temperature, very often observed for $\delta < \delta_{\max}$, can be explained by the high ratio U/V . It should be noted that in the considered case, the function $G(T)$ weakly depends on the temperature, which is also in agreement with the experimental data [66].

The very important feature of the considered approach is related to the existence of the non-zero value of the energy gap above the critical temperature. The performed numerical calculations show that $G(T)$ vanishes at the Nernst temperature (T^{**}) [47].

In the present paper, basing on knowledge of the experimental values of the critical temperature and the energy gap close to zero Kelvin, the full form of the function $G(T)$ has been determined. In particular, the following electron-doped superconductors, in the form of thin layers, have been taken into account: LCCO, $\text{Pr}_{2-x}\text{Ce}_x\text{CuO}_4$ (PCCO), and $\text{Nd}_{2-x}\text{Ce}_x\text{CuO}_4$ (NCCO). Additionally, it has been assumed that $\varepsilon_0 = 0$ and t_0 is equal to 380 meV [68–71].

2 Results

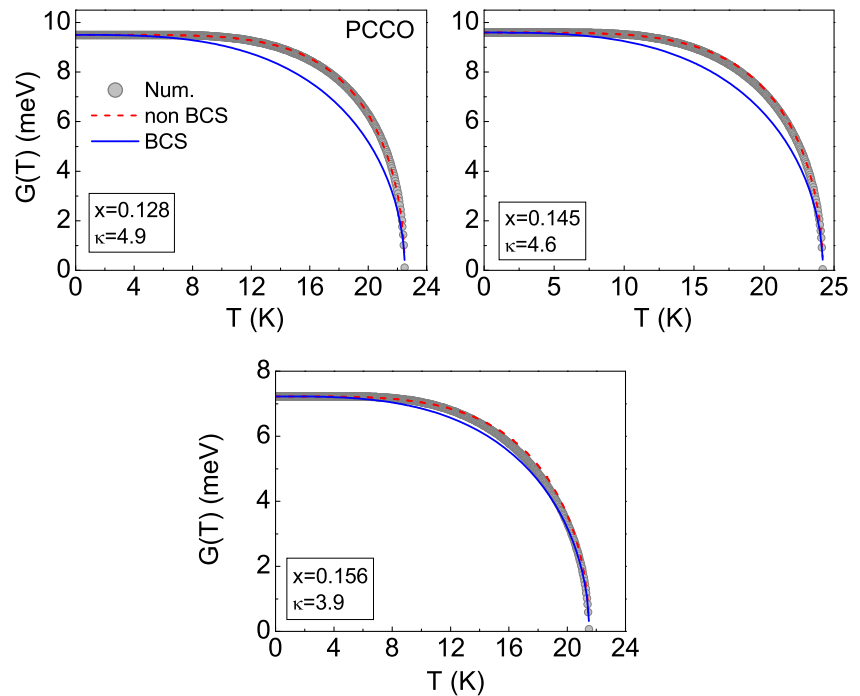
In Fig. 1a,b, the dependence of the critical temperature and the low-temperature value of the energy gap ($T = 1.6$ K) on the concentration of cerium has been presented. The symbols have been prepared on the basis of the experimental data [17, 67] (see also Table 1).

In the case of the critical temperature, it may be noted that the function $T_C(x)$ has the typical course observed in the high-temperature superconductors, corresponding approximately to the inverse parabola. The dependence of the energy gap on x is quite different—the value of $G(0)$ quickly decreases with the increase of the cerium concentration.

Table 2 The thermodynamic parameters of the PCCO superconductor

x	T_C (K)	$G(0)$ (meV)	V (meV)	$U^{(0)}$ (meV)	T^{**} (K)
0.128	22.5	9.5	380.4	51.5	—
0.145	24.2	9.6	391.6	42.2	—
0.156	21.5	7.2	373.6	12.0	—

Fig. 4 The dependence of the PCCO energy gap on the temperature for the selected values of the Ce concentration



Next, based on the data from the paper [17] and with the help of (6), the full dependence of the energy gap on the temperature has been calculated. Additionally, it has been assumed that the characteristic phonon frequency is equal to 29 meV [72]. The obtained results have been plotted in Fig. 2. It can be seen that for the lowest value for the concentration of cerium, the function $G(T)$ sharply differs from the BCS dependence of the energy gap on the temperature [34, 35]. In particular, G is much less dependent on the temperature and it does not vanish at the critical temperature, but at the Nernst temperature ($T^{**} = 28.9$ K). For the higher concentrations of cerium, there was no existence of

the Nernst area, although the courses of G on T also cannot be reproduced in the framework of the BCS model.

To demonstrate this, we note that the BCS relationship between the energy gap and the temperature can be described by the formula:

$$G(T) = G(0) \sqrt{1 - \left(\frac{T}{T_c}\right)^\kappa}, \quad (7)$$

where the exponent $[\kappa]_{\text{BCS}} = 3$ [73].

In the case of LCCO, the values of κ , which the best reproduce the numerical results, have been presented in

Fig. 5 The influence of the cerium concentration on the value of **a** the critical temperature and **b** the energy gap for the NCCO superconductor [67, 76]

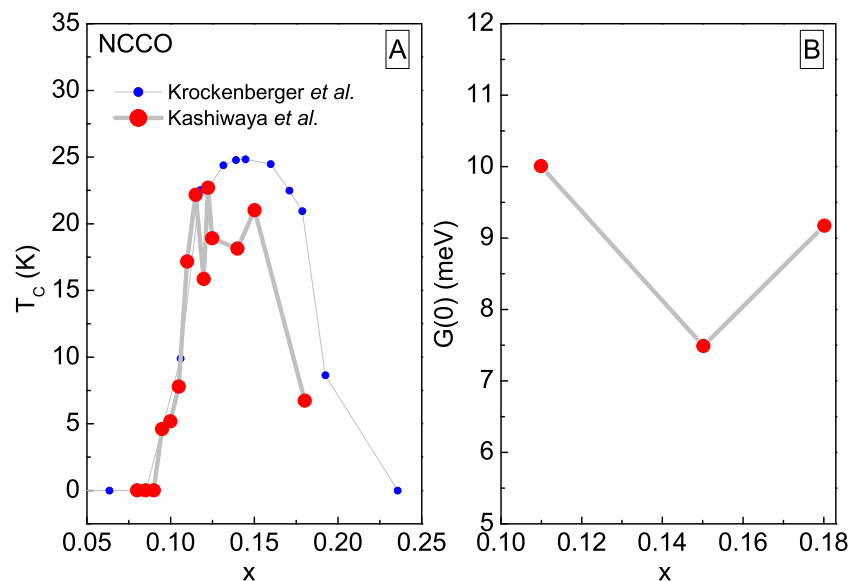


Table 3 The thermodynamic parameters of the NCCO superconductor

x	T_C (K)	$G(0)$ (meV)	V (meV)	U (meV)	T^{**} (K)
0.11	17.2	10.0	356.8	107.7	17.4
0.15	21.0	7.5	386.5	22.6	—
0.18	6.8	9.2	259.0	398.5	18.3

Fig. 2. The corresponding curves obtained by using (7) have been also plotted (the red dashed lines). The blue lines denote the BCS results. Additionally, let us pay attention that the values of κ , with the good approximation, can be calculated from: $\kappa \simeq G(0)/k_B T_C$.

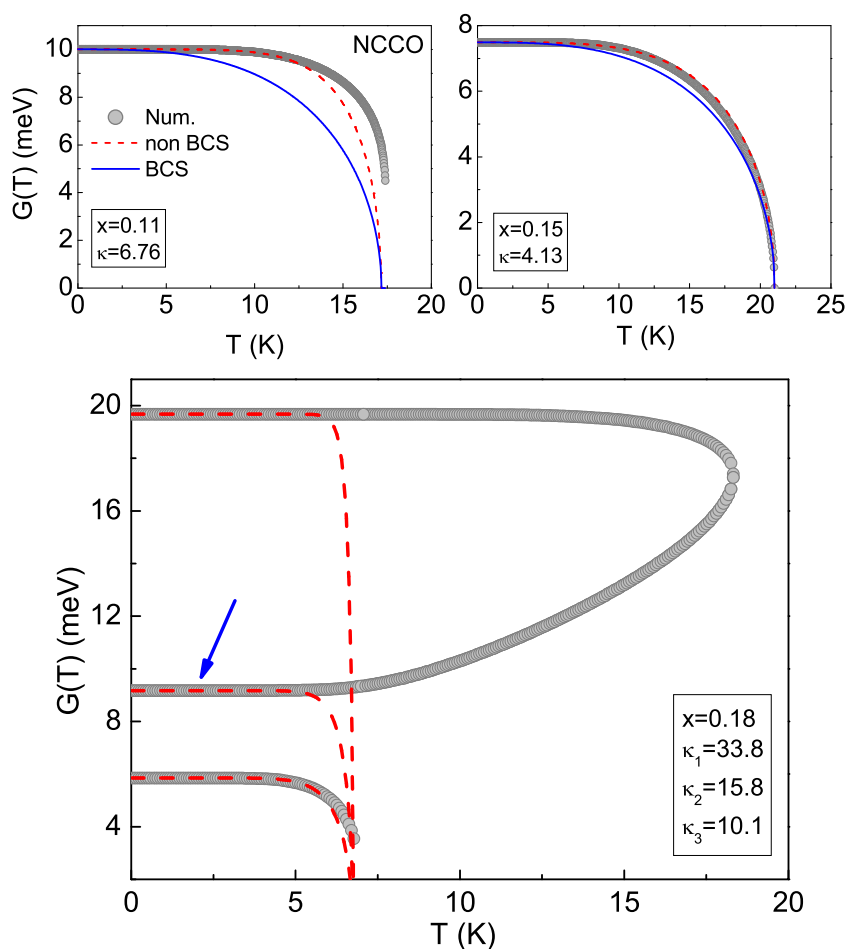
Analysing the data plotted in Fig. 2, it has been found that with the increasing concentration of cerium, the value of κ falls closer to the result of the BCS model. However, for the sufficiently high concentration ($x=0.135$), we have $\kappa < [\kappa]_{\text{BCS}}$. This is the interesting case, because it occurs when the effective potential $U^{(0)}$ changes its sign (see Table 1). From the physical point of view, the obtained result means that electron depairing correlations are stronger than

the phonon exchange processes in the EEPH channel ($U_0 > V_{U_0}$).

In the next step, the dependence of the energy gap on the temperature for the PCCO superconductor has been determined. The study has been based on the experimental results presented in Fig. 3 and in Table 2 (data from the paper of Skinta et al. [17]). During the calculations, it has been assumed that the maximum phonon frequency is equal to 33 meV [74, 75].

In the case of the PCCO superconductor, for the low concentrations of cerium, there was no Nernst phase. Additionally, the potential $U^{(0)}$ does not change the sign for any value of x . As it was in the case for the LCCO

Fig. 6 The dependence of the NCCO energy gap on the temperature for the selected values of the Ce concentration. The blue arrow indicates the branch of the energy gap, which is realized physically



superconductor, together with the increasing concentration of cerium, the exponent κ decreases and comes close to the value predicted by the BCS theory (see Fig. 4).

In the last step, the dependence of the energy gap on the temperature for the NCCO superconductor has been determined. The experimental results, on which the calculation has been based, are collected in Fig. 5 and in Table 3 (the paper of Kashiwaya et al. [76]). It has been adopted that the characteristic phonon frequency is equal to 28 meV [77].

The results presented in Fig. 6 prove that for the lowest considered concentration of cerium ($x=0.11$), above the critical temperature there is the Nernst region, which disappears for $T^{**} = 17.4$ K.

Very unusual results have been obtained for the highest concentration ($x=0.18$), which follows directly from the experimentally observed increase in the low-temperature energy gap value ($T = 4.5$ K), (see Fig. 5b or Table 3).

In this case, our model predicts that below the critical temperature there are three non-zero branches of $G(T)$, which vary widely in the adopted values (Fig. 6). Mathematically, it can be shown that the thermodynamic potential minimizes the branch with the highest values [47]. Experimental data demonstrate, however, that physically realized is the middle branch, which is due to the structural instability of the studied material (above $x>0.18$, it is impossible to synthesize the thin film of NCCO [76]).

In our opinion, the experimentally observed result is of fundamental importance in the process of falsification of the presented model. Namely, the obtained theoretical data clearly predict anomalous dependence of the energy gap on the temperature due to the fact of the increase in the value of G together with the increasing T .

In addition, note that above the superconducting state exists the wide Nernst region with the high value of the temperature T^{**} equal to 18.3 K. These predictions can be relatively easily checked experimentally, to which we strongly encourage.

Finally, it should be emphasized that for all concentrations of cerium, the NCCO exponent κ significantly differs from the value predicted by the BCS model.

3 Summary

In the paper, the full dependence of the energy gap on the temperature for the high-temperature superconductors LCCO, PCCO, and NCCO has been determined (the thin films). A wide range of the concentration of cerium has been taken into account.

It has been shown that generally, the curves $G(T)$ are the more different from the BCS function, than the lower concentration of cerium has been taken into consideration.

However, there are the derogation from the above scheme due to the structural instability of the system NCCO for $x = 0.18$. In this case, the increase of the energy gap together with the increasing of the temperature and the high value of T^{**} has been predicted.

For the superconductors LCCO and NCCO (the lowest concentration of cerium) induces the Nernst region, which disappears at the temperature slightly higher than T_C .

In the future, the thermodynamic properties of the electron-doped cuprates will be analysed in the framework of the extended Eliashberg approach [78]. The discussed procedure will enable the analysis of the strong-coupling and the retardation effects [79].

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